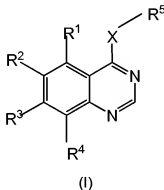


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

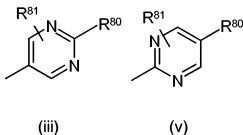
1. (Currently amended) A compound of formula (I)



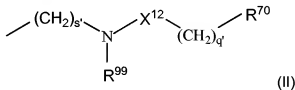
or a salt thereof;

where X is O, or S, S(O), S(O)₂ or NR⁶ where R⁶ is hydrogen or C₁₋₆alkyl;

R⁵ is [[or]] a group of sub-formula (iii) or (v)



where R⁸⁰ is a group of sub-formula (II)

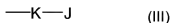


where q' is 0, 1, 2, 3 or 4;

s' is 0 or 1;

X¹² is C(O) or S(O₂),

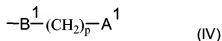
R⁷⁰ is hydrogen, hydroxy, C₁₋₆alkyl, C₁₋₆alkoxy, amino, N-C₁₋₆alkylamino, N,N-(C₁₋₆alkyl)₂amino, hydroxyC₂₋₆alkoxy, C₁₋₆alkoxyC₂₋₆alkoxy, aminoC₂₋₆alkoxy, N-C₁₋₆alkylaminoC₂₋₆alkoxy, N,N-(C₁₋₆alkyl)₂aminoC₂₋₆alkoxy or C₃₋₇cycloalkyl,
or R⁷⁰ is of the Formula (III):



wherein J is aryl, heteroaryl or heterocyclyl and K is a bond, oxy, imino, *N*-(C₁₋₆alkyl)imino, oxyC₁₋₆alkylene, iminoC₁₋₆alkylene, *N*-(C₁₋₆alkyl)iminoC₁₋₆alkylene, -NHC(O)-, -SO₂NH-, -NHSO₂- or -NHC(O)-C₁₋₆alkylene-,

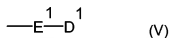
and any aryl, heteroaryl or heterocyclyl group in a R⁷⁰ group is optionally substituted by one or more groups selected from hydroxy, oxo, halo, trifluoromethyl, cyano, mercapto, nitro, amino, carboxy, carbamoyl, formyl, sulphamoyl, C₁₋₆alkyl, C₂₋₆alkenyl, C₂₋₆alkynyl, C₁₋₆alkoxy, -O-(C₁₋₃alkyl)-O-, C₁₋₆alkylS(O)_n- wherein n is 0-2, *N*-(C₁₋₆alkyl)amino, *N,N*-(C₁₋₆alkyl)₂amino, C₁₋₆alkoxycarbonyl, *N*-(C₁₋₆alkyl)carbamoyl, *N,N*-(C₁₋₆alkyl)₂carbamoyl, C₂₋₆alkanoyl, C₁₋₆alkanoyloxy, C₁₋₆alkanoylamino, *N*-(C₁₋₆alkyl)sulphamoyl, *N,N*-(C₁₋₆alkyl)₂sulphamoyl, C₁₋₆alkylsulphonylamino and C₁₋₆alkylsulphonyl-*N*-(C₁₋₆alkyl)amino,

or any aryl, heteroaryl or heterocyclyl group in a R⁷⁰ group is optionally substituted with one or more groups of the Formula (IV):



wherein A¹ is halo, hydroxy, C₁₋₆alkoxy, cyano, amino, *N*-(C₁₋₆alkyl)amino, *N,N*-(C₁₋₆alkyl)₂amino, carboxy, C₁₋₆alkoxycarbonyl, carbamoyl, *N*-(C₁₋₆alkyl)carbamoyl or *N,N*-(C₁₋₆alkyl)₂carbamoyl, p is 1 - 6, and B¹ is a bond, oxy, imino, *N*-(C₁₋₆alkyl)imino or -NHC(O)-, with the proviso that p is 2 or more unless B¹ is a bond or -NHC(O)-;

or any aryl, heteroaryl or heterocyclyl group in a R⁷⁰ group is optionally substituted with one or more groups of the Formula (V):



wherein D¹ is aryl, heteroaryl or heterocyclyl and E¹ is a bond, C₁₋₆alkylene, oxyC₁₋₆alkylene, oxy, imino, *N*-(C₁₋₆alkyl)imino, iminoC₁₋₆alkylene, *N*-(C₁₋₆alkyl)-iminoC₁₋₆alkylene, C₁₋₆alkylene-oxyC₁₋₆alkylene, C₁₋₆alkylene-iminoC₁₋₆alkylene, C₁₋₆alkylene-*N*-(C₁₋₆alkyl)-iminoC₁₋₆alkylene, -NHC(O)-, -NHSO₂-, -SO₂NH- or -NHC(O)-C₁₋₆alkylene-, and any aryl, heteroaryl or heterocyclyl group in a R⁷⁰ group is optionally substituted with one or more groups selected from hydroxy, halo, C₁₋₆alkyl, C₁₋₆alkoxy, carboxy, C₁₋₆alkoxycarbonyl, carbamoyl, *N*-(C₁₋₆alkyl)carbamoyl, *N*-(C₁₋₆alkyl)₂carbamoyl, C₂₋₆alkanoyl, amino, *N*-(C₁₋₆alkyl)amino and *N,N*-(C₁₋₆alkyl)₂amino, and any C₃₋₇cycloalkyl or heterocyclyl group in a R⁷⁰ group is optionally substituted with one or two oxo or thioxo substituents,

and any of the R⁷⁰ groups defined hereinbefore which comprises a CH₂ group which is attached to 2 carbon atoms or a CH₃ group which is attached to a carbon atom may optionally bear on each said CH₂ or CH₃ group a substituent selected from hydroxy, amino, C₁₋₆alkoxy,

N-C₁₋₆alkylamino, *N,N*-(C₁₋₆alkyl)₂amino and heterocyclyl;

or R⁷⁰ may be cycloalkenyl or alkenyl optionally substituted by aryl;

and R⁹⁹ is hydrogen or a group C(O)R⁷⁰ where R⁷⁰ is as defined above;

and

R⁸¹ is hydrogen, halo, C₁₋₄alkoxy, cyano, trifluoromethyl, or phenyl, and

R¹, R², R³, R⁴ are independently selected from halogeno, cyano, nitro, C₁₋₃alkylsulphanyl,

-N(OH)R⁷- wherein R⁷ is hydrogen, or C₁₋₃alkyl, or R⁹X¹- wherein X¹ represents a direct bond,

-O-, -CH₂-, -OC(O)-, -C(O)-, -S-, -SO-, -SO₂-, -NR¹⁰C(O)-, -C(O)NR¹¹-, -SO₂NR¹²-, -NR¹³SO₂- or

-NR¹⁴-, wherein R¹⁰, R¹¹, R¹², R¹³ and R¹⁴ each independently represents hydrogen, C₁₋₃alkyl or

C₁₋₃alkoxyC₂₋₃alkyl, provided that at least one of R¹, R², R³ and R⁴ is a group R⁹X¹- and R⁹ is

selected from one of the following groups: provided that at least one of R² or R³ is other than

hydrogen;

1) hydrogen or C₁₋₅alkyl which may be unsubstituted or which may be substituted with one or more groups selected from hydroxy, oxiranyl, fluoro, chloro, bromo and amino;

2) -R⁹X²C(O)R¹⁵ wherein X² represents -O- or -NR¹⁶- in which R¹⁶ represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl and R¹⁵ represents C₁₋₃alkyl, -NR¹⁷R¹⁸ or -OR¹⁹ wherein R¹⁷, R¹⁸ and R¹⁹ which may be the same or different each represents hydrogen, C₁₋₅alkyl, hydroxyC₁₋₅alkyl or C₁₋₃alkoxyC₂₋₃alkyl;

3) -R^bX³R²⁰ wherein X³ represents -O-, C(O)-S-, -SO-, -SO₂-, -OC(O)-, -NR²¹C(O)-, -C(O)NR²²-, -SO₂NR²³-, -NR²⁴SO₂- or -NR²⁵- wherein R²¹, R²², R²³, R²⁴ and R²⁵ each independently

represents hydrogen, C₁₋₃alkyl, hydroxy C₁₋₄alkyl or C₁₋₃alkoxyC₂₋₃alkyl and s is 1 or 2 and R²⁰

represents hydrogen, C₁₋₆alkyl, C₂₋₆alkenyl, cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl,

phenyl or a 5-6-membered saturated heterocyclic group with 1-2 heteroatoms, selected

independently from O, S and N, which C₁₋₆alkyl group may bear 1, 2 or 3 substituents selected

from oxo, hydroxy, halogeno, cyclopropyl, amino, C₁₋₄alkylamino, C₁₋₄alkanoyldi-C₁₋₄alkylamino,

C₁₋₄alkylthio, C₁₋₄alkoxy and which cyclic group may bear 1 or 2 substituents selected from oxo,

hydroxy, halogeno, cyano, C₁₋₄cyanoalkyl, C₁₋₄alkyl, C₁₋₄hydroxyalkyl, C₁₋₄alkoxy,

C₁₋₄alkoxyC₁₋₄alkyl, C₁₋₄alkylsulphonylC₁₋₄alkyl, C₁₋₄alkoxycarbonyl, C₁₋₄aminoalkyl,

C₁₋₄alkylamino, di(C₁₋₄alkyl)amino, C₁₋₄alkylaminoC₁₋₄alkyl, di(C₁₋₄alkyl)aminoC₁₋₄alkyl,

C₁₋₄alkylaminoC₁₋₄alkoxy, di(C₁₋₄alkyl)aminoC₁₋₄alkoxy and a group -(-O-)(R^b)_fD² wherein f is 0

or 1, g is 0 or 1 and D² is a C₃₋₆cycloalkyl group or a 5-6-membered saturated heterocyclic group

with 1-2 heteroatoms, selected independently from O, S and N, which cyclic group may bear one or more substituents selected from C₁₋₄alkyl;

4) -R^cX⁴R^{c'}X⁵R²⁶ wherein X⁴ and X⁵ which may be the same or different are each -O-, C(O), -S-, -SO-, -SO₂-, -NR²⁷C(O)_s-, -C(O)_sNR²⁸-, -SO₂NR²⁹-, -NR³⁰SO₂- or -NR³¹- wherein R²⁷, R²⁸, R²⁹, R³⁰ and R³¹ each independently represents hydrogen, C₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl and s is 1 or 2 and R²⁶ represents hydrogen, C₁₋₃alkyl, hydroxyC₁₋₃alkyl or C₁₋₃alkoxyC₂₋₃alkyl;

5) R³² wherein R³² is a 4-6-membered cycloalkyl or saturated heterocyclic ring, linked via carbon or nitrogen, with 1-2 heteroatoms, selected independently from O, S and N, which cycloalkyl or heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, cyano, C₁₋₄alkyl, hydroxyC₁₋₄alkyl, ~~C₁₋₄alkoxyC₁₋₄alkyl~~, cyanoC₁₋₄alkyl, cyclopropyl, C₁₋₄alkylsulphonylC₁₋₄alkyl, C₁₋₄alkoxycarbonyl, carboxamido, C₁₋₄aminoalkyl, C₁₋₄alkylamino, di(C₁₋₄alkyl)amino, C₁₋₄alkylaminoC₁₋₄alkyl, C₁₋₄alkanoyl, di(C₁₋₄alkyl)aminoC₁₋₄alkyl, C₁₋₄alkylaminoC₁₋₄alkoxy, di(C₁₋₄alkyl)aminoC₁₋₄alkoxy, nitro, amino, C₁₋₄alkoxy, C₁₋₄hydroxyalkoxy, carboxy, trifluoromethyl, -C(O)NR³⁸R³⁹, -NR⁴⁰C(O)R⁴¹, wherein R³⁸, R³⁹, R⁴⁰ and R⁴¹, which may be the same or different, each represents hydrogen, C₁₋₄alkyl, hydroxyC₁₋₄alkyl or C₁₋₃alkoxyC₂₋₃alkyl, and a group -(O-)_f(C₁₋₄alkyl)_gringD wherein f is 0 or 1, g is 0 or 1 and ring D is a cyclic group selected from C₃₋₆cycloalkyl, aryl or 5-6-membered saturated or unsaturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which cyclic group may bear one or more substituents selected from halo and C₁₋₄alkyl;

6) -R^cR³² wherein R³² is as defined hereinbefore;

7) -R^eR³² wherein R³² is as defined hereinbefore;

8) -R^fR³² wherein R³² is as defined hereinbefore;

9) R³³ wherein R³³ represents a pyridone group, a phenyl group or a 5-6-membered aromatic heterocyclic group, linked via carbon or nitrogen, with 1-3 heteroatoms selected from O, N and S, which pyridone, phenyl or aromatic heterocyclic group may carry up to 5 substituents selected from hydroxy, nitro, halogeno, amino, C₁₋₄alkyl, C₁₋₄alkoxy, C₁₋₄hydroxyalkyl, C₁₋₄aminoalkyl, C₁₋₄alkylamino, C₁₋₄hydroxyalkoxy, oxo, cyanoC₁₋₄alkyl, cyclopropyl, C₁₋₄alkylsulphonylC₁₋₄alkyl, C₁₋₄alkoxycarbonyl, di(C₁₋₄alkyl)amino, C₁₋₄alkylaminoC₁₋₄alkyl, C₁₋₄alkanoyl, di(C₁₋₄alkyl)aminoC₁₋₄alkyl, C₁₋₄alkylaminoC₁₋₄alkoxy, di(C₁₋₄alkyl)aminoC₁₋₄alkoxy, carboxy, carboxamido, trifluoromethyl, cyano, -C(O)NR³⁸R³⁹, -NR⁴⁰C(O)R⁴¹, wherein R³⁸, R³⁹, R⁴⁰ and R⁴¹, which may be the same or different, each represents hydrogen, C₁₋₄alkyl, hydroxyC₁₋₄alkyl or C₁₋₃alkoxyC₂₋₃alkyl, and a group -(O-)_f(C₁₋₄alkyl)_gringD wherein f is 0 or 1, g is 0 or 1 and ring D is a cyclic group selected from C₃₋₆cycloalkyl, aryl or 5-6-membered saturated or unsaturated

heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which cyclic group may bear one or more substituents selected from halo and $C_{1-4}alkyl$;

10) $-R^gR^{33}$ wherein R^{33} is as defined hereinbefore;

11) $-R^hR^{33}$ wherein R^{33} is as defined hereinbefore;

12) $-R^iR^{33}$ wherein R^{33} is as defined hereinbefore;

13) $-R^jX^6R^{33}$ wherein X^6 represents $-O-$, $-C(O)-$, $-S-$, $-SO-$, $-SO_2-$, $-OC(O)-$, $-NR^{38'}C(O)-$, $-C(O)NR^{39'}$, $-SO_2NR^{40'}$, $-NR^{41'}SO_2-$ or $-NR^{42'}$, wherein $R^{38'}$, $R^{39'}$, $R^{40'}$, $R^{41'}$ and $R^{42'}$ each independently represents hydrogen, $C_{1-3}alkyl$, hydroxy $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$, and R^{33} is as defined hereinbefore;

14) $-R^kX^7R^{33}$ wherein X^7 represents $-O-$, $C(O)-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{43}C(O)-$, $-C(O)NR^{44-}$, $-SO_2NR^{45-}$, $-NR^{46}SO_2-$ or $-NR^{47-}$, wherein R^{43} , R^{44} , R^{45} , R^{46} and R^{47} each independently represents hydrogen, $C_{1-3}alkyl$, hydroxy $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$, and R^{33} is as defined hereinbefore;

15) $-R^mX^8R^{33}$ wherein X^8 represents $-O-$, $-C(O)-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{48}C(O)-$, $-C(O)NR^{49-}$, $-SO_2NR^{50-}$, $-NR^{51}SO_2-$ or $-NR^{52-}$, wherein R^{48} , R^{49} , R^{50} , R^{51} and R^{52} each independently represents hydrogen, $C_{1-3}alkyl$, hydroxy $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$, and R^{33} is as defined hereinbefore;

16) $-R^nX^9R^{33}$ wherein X^9 represents $-O-$, $-C(O)-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{53}C(O)-$, $-C(O)NR^{54-}$, $-SO_2NR^{55-}$, $-NR^{56}SO_2-$ or $-NR^{57-}$, wherein R^{53} , R^{54} , R^{55} , R^{56} and R^{57} each independently represents hydrogen, $C_{1-3}alkyl$, hydroxy $C_{1-3}alkyl$ or $C_{1-3}alkoxyC_{2-3}alkyl$, and R^{33} is as defined hereinbefore;

17) $-R^pX^9-R^{32}$ wherein X^9 and R^{32} are as defined hereinbefore;

18) $C_{2-5}alkenyl$ which may be unsubstituted or which may be substituted with one or more groups selected from hydroxy, fluoro, amino, $C_{1-4}alkylamino$, $N,N-di(C_{1-4}alkyl)amino$, aminosulphonyl, $N-C_{1-4}alkylaminosulphonyl$ and $N,N-di(C_{1-4}alkyl)aminosulphonyl$;

19) $C_{2-5}alkynyl$ which may be unsubstituted or which may be substituted with one or more groups selected from hydroxy, fluoro, amino, $C_{1-4}alkylamino$, $N,N-di(C_{1-4}alkyl)amino$, aminosulphonyl, $N-C_{1-4}alkylaminosulphonyl$ and $N,N-di(C_{1-4}alkyl)aminosulphonyl$;

20) $-R^qX^9R^{32}$ wherein X^9 and R^{32} are as defined hereinbefore;

21) $-R^qX^9R^qR^{32}$ wherein X^9 and R^{32} are as defined hereinbefore; and

22) $-R^qR^{58}(R^q)_q(X^9)$, R^{59} wherein X^9 is as defined hereinbefore, q is 0 or 1, r is 0 or 1, and R^{58} is a $C_{1-3}alkylene$ group or a cyclic group selected from cyclopropyl, cyclobutyl, cyclopentylene, cyclohexylene or a 5-6-membered saturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which $C_{1-3}alkylene$ group may bear 1 or 2 substituents selected

from oxo, hydroxy, halogeno and C₁₋₄alkoxy and which cyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, cyano, C₁₋₄cyanoalkyl, C₁₋₄alkyl, C₁₋₄hydroxyalkyl, C₁₋₄alkoxy, C₁₋₄alkoxyC₁₋₄alkyl, C₁₋₄alkylsulphonylC₁₋₄alkyl, C₁₋₄alkoxycarbonyl, C₁₋₄aminoalkyl, C₁₋₄alkylamino, di(C₁₋₄alkyl)amino, C₁₋₄alkylaminoC₁₋₄alkyl, di(C₁₋₄alkyl)aminoC₁₋₄alkyl, C₁₋₄alkylaminoC₁₋₄alkoxy, di(C₁₋₄alkyl)aminoC₁₋₄alkoxy and a group $-(O)_f(C_{1-4}alkyl)_gringD$, wherein f is 0 or 1, g is 0 or 1 and ring D is a cyclic group selected from C₃₋₆cycloalkyl, aryl or 5-6-membered saturated or unsaturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which cyclic group may bear one or more substituents selected from halo and C₁₋₄alkyl; and R⁵⁹ is hydrogen, C₁₋₃alkyl, or a cyclic group selected from cyclopropyl, cyclobutyl, cyclopentyl, cyclohexyl and a 5-6-membered saturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which C₁₋₃alkyl group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C₁₋₄alkoxy and which cyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, cyano, C₁₋₄cyanoalkyl, C₁₋₄alkyl, C₁₋₄hydroxyalkyl, C₁₋₄alkoxy, C₁₋₄alkoxyC₁₋₄alkyl, C₁₋₄alkylsulphonylC₁₋₄alkyl, C₁₋₄alkoxycarbonyl, C₁₋₄aminoalkyl, C₁₋₄alkylamino, di(C₁₋₄alkyl)amino, C₁₋₄alkylaminoC₁₋₄alkyl, di(C₁₋₄alkyl)aminoC₁₋₄alkyl, C₁₋₄alkylaminoC₁₋₄alkoxy, di(C₁₋₄alkyl)aminoC₁₋₄alkoxy and a group $-(O)_f(C_{1-4}alkyl)_gringD$ wherein f is 0 or 1, g is 0 or 1 and D is a cyclic group selected from C₃₋₆cycloalkyl, aryl or 5-6-membered saturated or unsaturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which cyclic group may bear one or more substituents selected from halo and C₁₋₄alkyl;

and wherein R^a, R^b, R^{b'}, R^c, R^{c'}, R^d, R^g, Rⁱ, Rⁿ, R^p, R^{p'}, R^r, R^u, R^v and R^{v'} are independently selected from C₁₋₈alkylene groups optionally substituted by one or more substituents selected from hydroxy, halogeno, amino,

R^e, R^h, R^k and R^l are independently selected from C₂₋₈alkenylene groups optionally substituted by one or more substituents selected from hydroxy, halogeno, amino, and R^l may additionally be a bond; and

R^f, Rⁱ, R^m and R^u are independently selected from C₂₋₈alkynylene groups optionally substituted by one or more substituents selected from hydroxy, halogeno, amino; and

where a functional group is selected from nitro, cyano, halo, oxo, =CR⁷⁸R⁷⁹, C(O),R⁷⁷, OR⁷⁷, S(O),R⁷⁷, NR⁷⁸R⁷⁹, C(O)NR⁷⁸R⁷⁹, OC(O)NR⁷⁸R⁷⁹, =NOR⁷⁷, -NR⁷⁷C(O),R⁷⁸, -NR⁷⁷CONR⁷⁸R⁷⁹, -N=CR⁷⁸R⁷⁹, S(O),NR⁷⁸R⁷⁹ or -NR⁷⁷S(O),R⁷⁸ where R⁷⁷, R⁷⁸ and R⁷⁹ are independently selected from hydrogen, optionally substituted hydrocarbyl, optionally substituted heterocyclyl or optionally substituted alkoxy, or R⁷⁸ and R⁷⁹ together form an optionally substituted ring which optionally contains further heteroatoms such as oxygen, nitrogen, S, S(O) or S(O)₂, where x is

an integer of 1 or 2, y is 0 or an integer of 1-3 and where hydrocarbonyl, heterocyclyl or alkoxy groups R^{77} , R^{78} and R^{79} as well as rings formed by R^{78} and R^{79} are optionally substituted by halo, perhaloalkyl, mercapto, alkylthio, hydroxy, carboxy, alkoxy, heteroaryl, heteroaryloxy, cycloalkyl, cycloalkenyl, cycloalkynyl, alkenyloxy, alkynyloxy, alkoxyalkoxy, aryloxy where the aryl group may be substituted by halo, nitro, or hydroxy, cyano, nitro, amino, mono- or di-alkyl amino, oximino or $S(O)_yR^{90}$ where y is 0 or an integer of 1-3 and R^{90} is a alkyl; and wherein hydrocarbonyl is selected from alkyl, alkenyl, alkynyl, aryl, aralkyl, cycloalkyl, cycloalkenyl, or combinations thereof.

2-5. (Canceled)

6. (Currently amended) A compound according to claim 1 wherein R^1 , R^2 , R^3 , R^4 are independently selected from, halo, cyano, nitro, trifluoromethyl, C_{1-3} alkyl, or other groups from formula $-X^1R^9$ wherein X^1 represents a direct bond, $-O-$, $-CH_2-$, $-OCO-$, carbonyl, $-S-$, $-SO-$, $-SO_2-$, $-NR^{10}CO-$, $-CONR^{11}-$, $-SO_2NR^{12}-$, $-NR^{13}SO_2-$ or $-NR^{14}-$, wherein R^{10} , R^{11} , R^{12} , R^{13} and R^{14} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl, and R^9 is selected from one of the following groups:

1') hydrogen or C_{1-5} alkyl which may be unsubstituted or which may be substituted with one or more groups selected from hydroxy, fluoro or amino,

2') C_{1-5} alkyl $X^2C(O)R^{15}$ wherein X^2 represents $-O-$ or $-NR^{16}-$ in which R^{15} represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl and R^{16} represents C_{1-3} alkyl, $-NR^{17}R^{18}$ or $-OR^{19}$ wherein R^{17} , R^{18} and R^{19} which may be the same or different each represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl;

3') C_{1-5} alkyl X^3R^{20} wherein X^3 represents $-O-$, $-S-$, $-SO-$, $-SO_2-$, $-OCO-$, $-NR^{21}CO-$, $-CONR^{22}-$, $-SO_2NR^{23}-$, $-NR^{24}SO_2-$ or $-NR^{25}-$, wherein R^{21} , R^{22} , R^{23} , R^{24} and R^{25} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl and R^{20} represents hydrogen, C_{1-3} alkyl, cyclopentyl, cyclohexyl or a 5-6-membered saturated heterocyclic group with 1-2 heteroatoms, selected independently from O, S and N, which C_{1-3} alkyl group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno and C_{1-4} alkoxy and which cyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C_{1-4} alkyl, C_{1-4} hydroxyalkyl and C_{1-4} alkoxy;

4') C_{1-5} alkyl X^4C_{1-5} alkyl X^5R^{26} wherein X^4 and X^5 which may be the same or different are each $-O-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{27}CO-$, $-CONR^{28}-$, $-SO_2NR^{29}-$, $-NR^{30}SO_2-$ or $-NR^{31}-$, wherein R^{27} , R^{28} , R^{29} , R^{30} and R^{31} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl and R^{26} represents hydrogen or C_{1-3} alkyl;

5') R^{32} wherein R^{32} is a 5-6-membered saturated heterocyclic group, linked via carbon or nitrogen, with 1-2 heteroatoms, selected independently from O, S and N, which heterocyclic group may bear 1 or 2 substituents selected from oxo, hydroxy, halogeno, C_{1-4} alkyl, C_{1-4} hydroxyalkyl, C_{1-4} alkoxy, C_{1-4} alkoxy C_{1-4} alkyl and C_{1-4} alkylsulphonyl C_{1-4} alkyl;

6') C_{1-5} alkyl R^{32} wherein R^{32} is as defined in (5') above;

7') C_{2-5} alkenyl R^{32} wherein R^{32} is as defined in (5') above;

8') C_{2-5} alkynyl R^{32} wherein R^{32} is as defined in (5') above;

9') R^{33} wherein R^{33} represents a pyridone group, a phenyl group or a 5-6-membered aromatic heterocyclic group, linked via carbon or nitrogen, with 1-3 heteroatoms selected from O, N and S, which pyridone, phenyl or aromatic heterocyclic group may carry up to 5 substituents on an available carbon atom selected from hydroxy, halogeno, amino, C_{1-4} alkyl, C_{1-4} alkoxy, C_{1-4} hydroxyalkyl, C_{1-4} aminoalkyl, C_{1-4} alkylamino, C_{1-4} hydroxyalkoxy, carboxy, trifluoromethyl, cyano, $-CONR^{384}R^{3936}$ and $-NR^{4036}COR^{4137}$, wherein R^{384} , R^{3936} , R^{4036} and R^{4137} , which may be the same or different, each represents hydrogen, C_{1-4} alkyl or C_{1-3} alkoxy C_{2-3} alkyl;

10') C_{1-5} alkyl R^{33} wherein R^{33} is as defined in (9') above;

11') C_{2-5} alkenyl R^{33} wherein R^{33} is as defined in (9') above;

12') C_{2-5} alkynyl R^{33} wherein R^{33} is as defined in (9') above;

13') C_{1-5} alkyl X^6R^{33} wherein X^6 represents $-O-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{38}CO-$, $-CONR^{39}-$, $-SO_2NR^{40}-$, $-NR^{41}SO_2-$ or $-NR^{42}-$, wherein R^{38} , R^{39} , R^{40} , R^{41} and R^{42} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl, and R^{33} is as defined hereinbefore;

14') C_{2-5} alkenyl X^7R^{33} wherein X^7 represents $-O-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{43}CO-$, $-CONR^{44}-$, $-SO_2NR^{45}-$, $-NR^{46}SO_2-$ or $-NR^{47}-$, wherein R^{43} , R^{44} , R^{45} , R^{46} and R^{47} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl, and R^{33} is as defined hereinbefore;

15') C_{2-5} alkynyl X^8R^{33} wherein X^8 represents $-O-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{48}CO-$, $-C(O)NR^{49}-$, $-SO_2NR^{50}-$, $-NR^{51}SO_2-$ or $-NR^{52}-$, wherein R^{48} , R^{49} , R^{50} , R^{51} and R^{52} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl, and R^{33} is as defined hereinbefore;

16') C_{1-3} alkyl X^9C_{1-3} alkyl R^{33} wherein X^9 represents $-O-$, $-S-$, $-SO-$, $-SO_2-$, $-NR^{53}CO-$, $-C(O)NR^{54}-$, $-SO_2NR^{55}-$, $-NR^{56}SO_2-$ or $-NR^{57}-$, wherein R^{53} , R^{54} , R^{55} , R^{56} and R^{57} each independently represents hydrogen, C_{1-3} alkyl or C_{1-3} alkoxy C_{2-3} alkyl, and R^{33} is as defined hereinbefore; and

17') C_{1-3} alkyl X^9C_{1-3} alkyl R^{32} wherein X^9 and R^{32} are as defined in (5') above, provided that at least one of R^2 or R^3 is other than hydrogen.

7. (Previously presented) A compound according to claim 1, where R^1 is hydrogen and R^4 is hydrogen, halo, C_{1-4} alkyl or C_{1-4} alkoxy.

8-9. (Canceled)

10. (Previously presented) A compound according to claim 1 or claim 7 wherein R^3 is a group X^1R^9 where X^1 is oxygen.

11. (Cancelled)

12. (Previously presented) A compound according to claim 7 wherein R^9 is selected from a group (1), (3), (6) or (10).

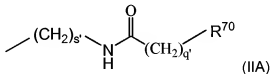
13. (Previously presented) A compound according to claim 12 wherein X is NH or O.

14-17. (Canceled)

18. (Previously presented) A compound according to claim 13 wherein R^5 is a group of formula (iii).

19-20. (Canceled)

21. (Currently amended) A compound according to claim 13 wherein R^{80} is a group of sub formula (II) which is a group of formula (IIA)



where s' , q' and R^{70} are as defined in claim 1,

where q' is 0, 1, 2, 3 or 4;

s' is 0 or 1;

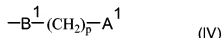
R^{70} is C_{3-7} cycloalkyl,

or R^{70} is of the Formula (III);



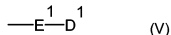
wherein J is aryl, heteroaryl or heterocyclyl and K is a bond, oxy, imino, $N-(C_{1-6}alkyl)imino$, $oxyC_{1-6}alkylene$, $iminoC_{1-6}alkylene$, $N-(C_{1-6}alkyl)iminoC_{1-6}alkylene$, $-NHC(O)-$, $-SO_2NH-$, $-NHSO_2-$ or $-NHC(O)-C_{1-6}alkylene-$,

and any aryl, heteroaryl or heterocyclyl group in a R^{70} group is optionally substituted by one or more groups selected from hydroxy, oxo, halo, trifluoromethyl, cyano, mercapto, nitro, amino, carboxy, carbamoyl, formyl, sulphamoyl, $C_{1-6}alkyl$, $C_{2-6}alkenyl$, $C_{2-6}alkynyl$, $C_{1-6}alkoxy$, $-O-(C_{1-3}alkyl)-O-$, $C_{1-6}alkylS(O)_n$ wherein n is 0-2, $N-C_{1-6}alkylamino$, $N,N-(C_{1-6}alkyl)_2amino$, $C_{1-6}alkoxycarbonyl$, $N-C_{1-6}alkylcarbamoyl$, $N,N-(C_{1-6}alkyl)_2carbamoyl$, $C_{2-6}alkanoyl$, $C_{1-6}alkanoyloxy$, $C_{1-6}alkanoylamino$, $N-C_{1-6}alkylsulphamoyl$, $N,N-(C_{1-6}alkyl)_2sulphamoyl$, $C_{1-6}alkylsulphonylamino$ and $C_{1-6}alkylsulphonyl-N-(C_{1-6}alkyl)amino$,
or any aryl, heteroaryl or heterocyclyl group in a R^{70} group is optionally substituted with one or more groups of the Formula (IV):



wherein A^1 is halo, hydroxy, $C_{1-6}alkoxy$, cyano, amino, $N-C_{1-6}alkylamino$, $N,N-(C_{1-6}alkyl)_2amino$, carboxy, $C_{1-6}alkoxycarbonyl$, carbamoyl, $N-C_{1-6}alkylcarbamoyl$ or $N,N-(C_{1-6}alkyl)_2carbamoyl$, p is 1 - 6, and B^1 is a bond, oxy, imino, $N-(C_{1-6}alkyl)imino$ or $-NHC(O)-$, with the proviso that p is 2 or more unless B^1 is a bond or $-NHC(O)-$;

or any aryl, heteroaryl or heterocyclyl group in a R^{70} group is optionally substituted with one or more groups of the Formula (V):



wherein D^1 is aryl, heteroaryl or heterocyclyl and E^1 is a bond, $C_{1-6}alkylene$, $oxyC_{1-6}alkylene$, oxy, imino, $N-(C_{1-6}alkyl)imino$, $iminoC_{1-6}alkylene$, $N-(C_{1-6}alkyl)iminoC_{1-6}alkylene$, $C_{1-6}alkyleneoxyC_{1-6}alkylene$, $C_{1-6}alkyleneiminoC_{1-6}alkylene$, $C_{1-6}alkylene-N-(C_{1-6}alkyl)iminoC_{1-6}alkylene$, $-NHC(O)-$, $-NHSO_2-$, $-SO_2NH-$ or $-NHC(O)-C_{1-6}alkylene-$, and any aryl, heteroaryl or heterocyclyl group in a R^{70} group is optionally substituted with one or more groups selected from hydroxy, halo, $C_{1-6}alkyl$, $C_{1-6}alkoxy$, carboxy, $C_{1-6}alkoxycarbonyl$, carbamoyl, $N-C_{1-6}alkylcarbamoyl$, $N-(C_{1-6}alkyl)_2carbamoyl$, $C_{2-6}alkanoyl$, amino, $N-C_{1-6}alkylamino$ and $N,N-(C_{1-6}alkyl)_2amino$,
and any $C_{3-7}cycloalkyl$ or heterocyclyl group in a R^{70} group is optionally substituted with one or two oxo or thioxo substituents,

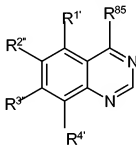
and any of the R^{70} groups defined hereinbefore which comprises a CH_2 group which is attached to 2 carbon atoms or a CH_3 group which is attached to a carbon atom may optionally bear on

each said CH₂ or CH₃ group a substituent selected from hydroxy, amino, C₁₋₆alkoxy, N-C₁₋₆alkylamino, N,N-(C₁₋₆alkyl)₂amino and heterocyclchl;
or R⁷⁰ may be cycloalkenyl.

22. (Previously presented) A compound according to claim 1 or claim 21 wherein R⁸⁰ includes a group R⁷⁰ and said group is phenyl optionally substituted by halo.

23-25. (Cancelled)

26. (Withdrawn) A method for preparing a compound of formula (I) as defined in claim 1, which method comprises reacting a compound of formula (VII)



(VII)

where R^{1'}, R^{2''}, R³, and R^{4'} are equivalent to a group R¹, R², R³ and R⁴ as defined in relation to formula (I), and R⁸⁵ is a leaving group, with a compound of formula (VIII)



where X and R⁵ are as defined in relation to formula (I).

27- 28. (Canceled)

29. (Currently amended) A pharmaceutical composition comprising a compound according to any one of claims 1, 7, 12, 18, 21 or 32~~34-6~~ or salt thereof, in combination with a pharmaceutically acceptable carrier.

30. (Canceled)

31. (Previously presented) A compound according to claim 1 wherein both R¹ and R⁴ are hydrogen.

32. (Previously presented) A compound according to claim 12 wherein one of R^2 or R^3 is 3-morpholinopropoxy.

33-36. (Cancelled)

37. (Previously presented) A method for treating colorectal or breast cancer in a warm blooded animal, such as man, in need of such treatment, which comprises administering to said animal an effective amount of a compound according to claim 1, or salt thereof.